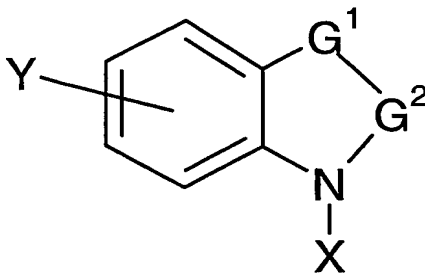


## WHAT IS CLAIMED IS:

1. A compound structurally represented by Formula I

5



(I)

or pharmaceutically acceptable salts thereof wherein:

$G^1$  is  $-\text{CH}_2-$ , or  $-\text{CH}_2\text{-CH}_2-$ ,

10  $G^2$  is  $-\text{CH}_2-$ , or  $-\text{C}(\text{O})-$ ,

or  $G^1$  and  $G^2$  taken together combine to form  $-\text{CH}=\text{CH}-$  or  $-\text{CH}_2\text{-CH}=\text{CH}-$ ,

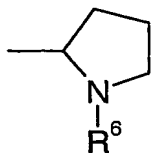
Y is

- 15  $-\text{OCH}_2\text{CH}_2\text{N-piperidinyl}$ ,  
 $-\text{OCH}_2\text{CH}_2\text{CH}_2\text{N-piperidinyl}$ ,  
 $-\text{OCH}_2\text{CH}_2\text{N-pyrrolidinyl}$ ,  
 $-\text{OCH}_2\text{CH}_2\text{CH}_2\text{N-pyrrolidinyl}$ ,

X is H,  $-\text{COR}^3$ ,  $-\text{CH}_2\text{R}^4$ ,  $-\text{SO}_2\text{R}^5$ ,

$\text{R}^3$  is

- 20  $-(\text{C}_1\text{-C}_8)$  alkyl, optionally substituted with 1 to 3 halogens,  
 $-(\text{C}_3\text{-C}_8)$  cycloalkyl, optionally substituted with 1 to 3 halogens,  
 $-\text{O}(\text{C}_1\text{-C}_8)$  alkyl, optionally substituted with 1 to 3 halogens,



, wherein  $R^6$  is  $-(C_1-C_6)$  alkyl, or  $-COO-(C_1-C_6)$  alkyl,

-Furanyl,

-Thienyl,

-NH-phenyl,

5 -NH- $(C_1-C_4)$ alkyl-phenyl,

-NH- $(C_1-C_8)$  alkyl, optionally substituted with 1 to 4 halogens,

-NH- $(C_3-C_8)$  cycloalkyl, optionally substituted once or twice with halogens,

-CH<sub>2</sub>-Pyridinyl,

-CH<sub>2</sub>N  $(C_1-C_6)$  alkyl  $(C_1-C_6)$  alkyl,

10 -CH<sub>2</sub>N-phenyl,

$R^4$  is

- $(C_1-C_8)$  alkyl, optionally substituted with 1 to 4 halogens,

- $(C_3-C_8)$  cycloalkyl,

- $(C_1-C_8)$  alkyl-NH<sub>2</sub>,

15 - $(C_1-C_4)$  alkyl -phenyl,

-CH<sub>2</sub>N-phenyl,

-phenyl-O- $(C_1-C_4)$  alkyl -phenyl,

- $(C_1-C_4)$  alkyl-O- $(C_1-C_4)$  alkyl-phenyl,

-CH<sub>2</sub>NCO<sub>2</sub>- $(C_1-C_6)$  alkyl,

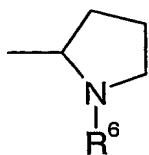
20 -Phenyl,

-Thienyl,

-Furanyl,

-Imidazolyl,

-Naphthyl,



, wherein R<sup>6</sup> is -(C<sub>1</sub>-C<sub>6</sub>) alkyl, or -COO-(C<sub>1</sub>-C<sub>6</sub>) alkyl,

-Biphenyl, and

R<sup>5</sup> is

-Phenyl,

5

-(C<sub>1</sub>-C<sub>4</sub>) alkyl,

-(C<sub>1</sub>-C<sub>4</sub>) alkyl -phenyl.

2. The compound of claim 1, wherein G1 is -CH<sub>2</sub>-.

3. The compound of claim 1, wherein G1 -CH<sub>2</sub>-CH<sub>2</sub>-.

4. The compound of claim 2, wherein Y is in the 5 position.

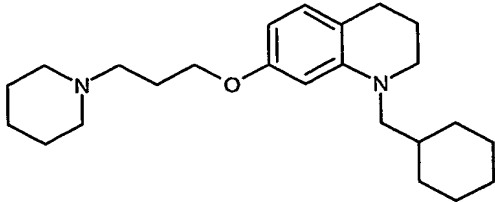
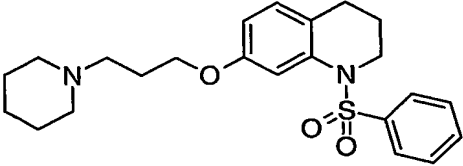
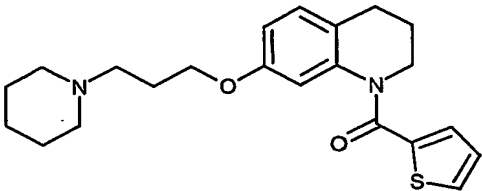
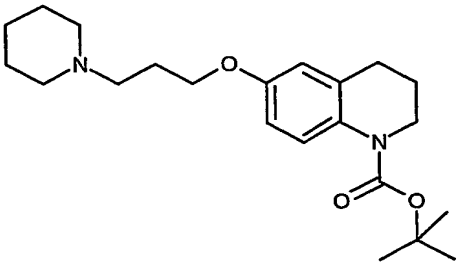
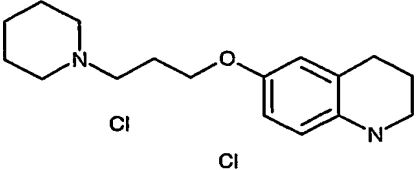
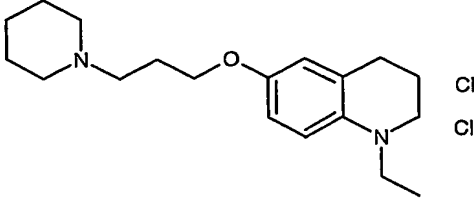
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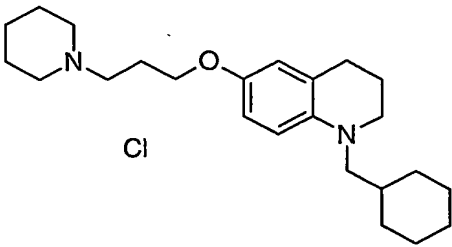
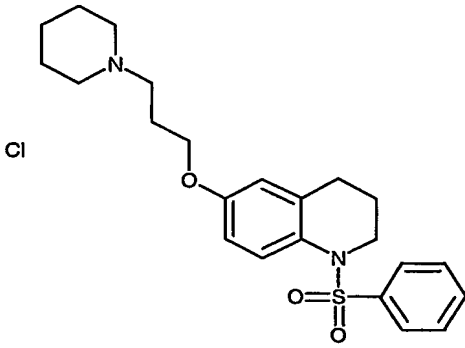
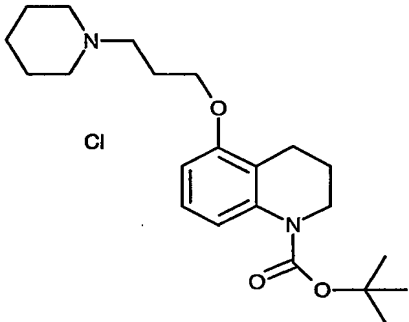
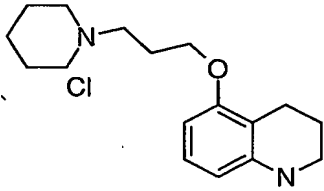
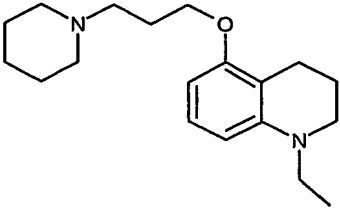
5. The compound of claim 3, wherein Y is in the 6 position.

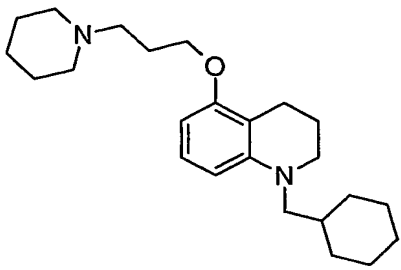
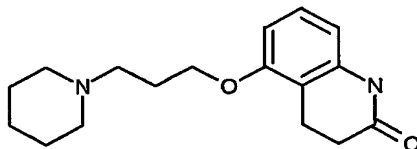
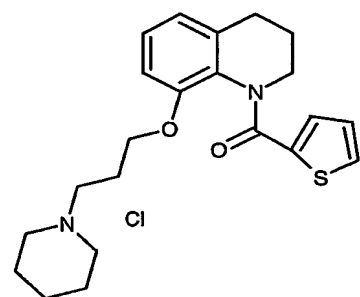
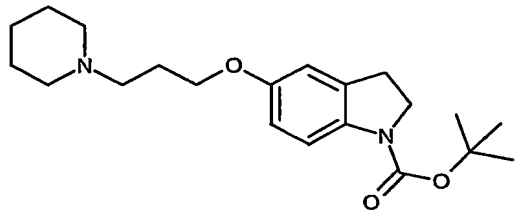
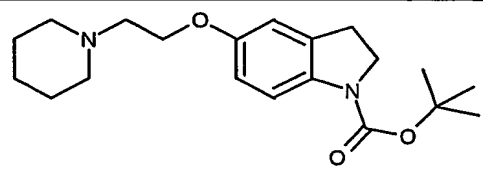
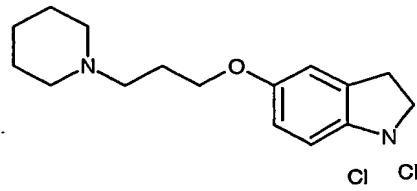
6. The compound of claim 4, wherein X is -COR<sup>3</sup>.

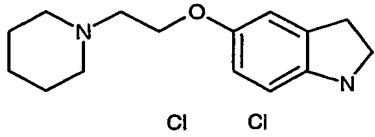
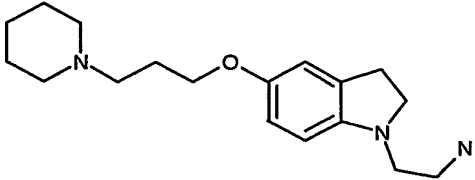
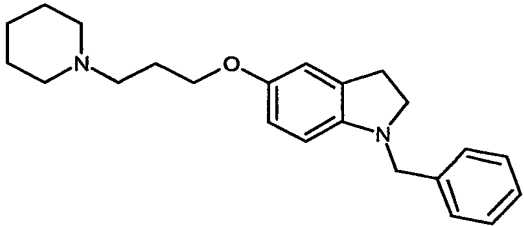
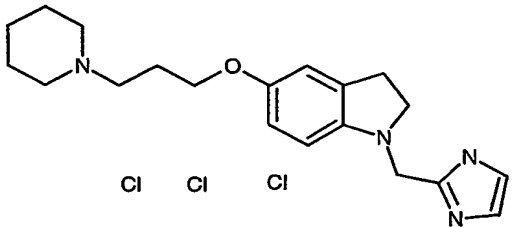
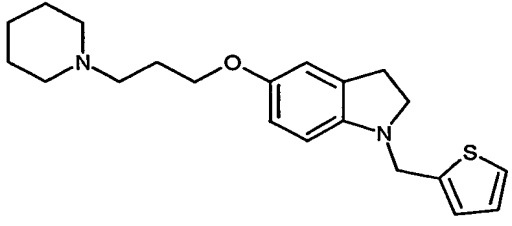
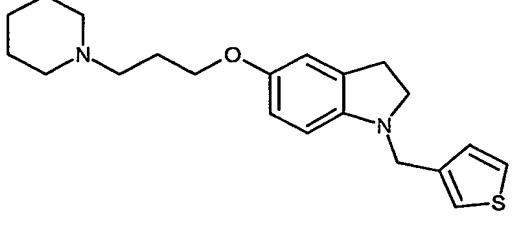
7. The compound of claim 1, selected from the group consisting of:

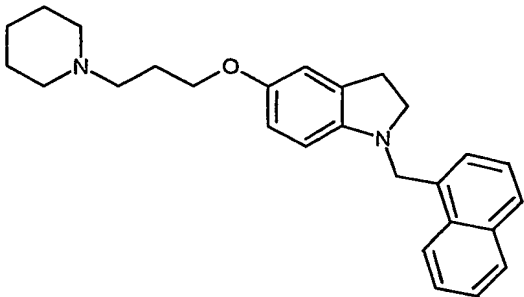
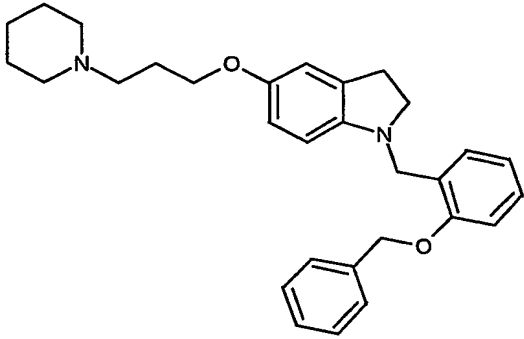
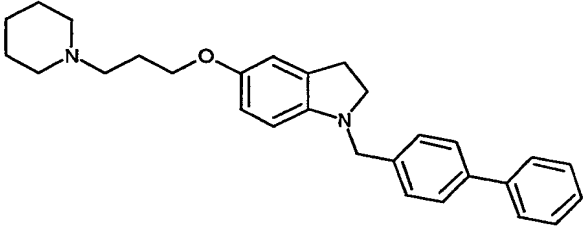
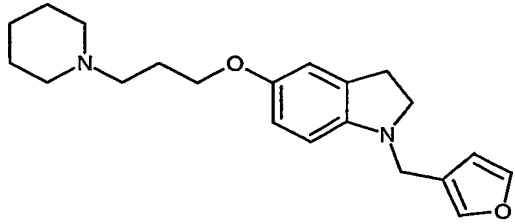
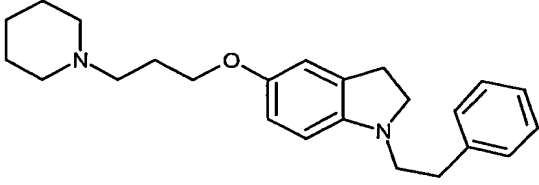
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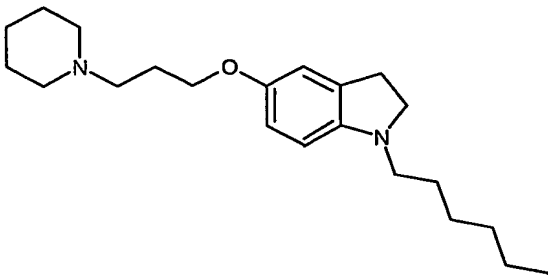
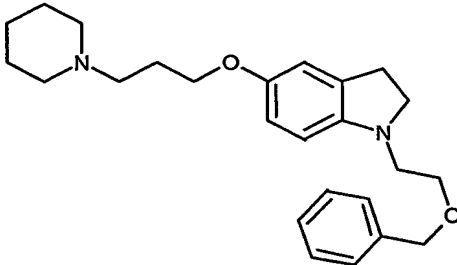
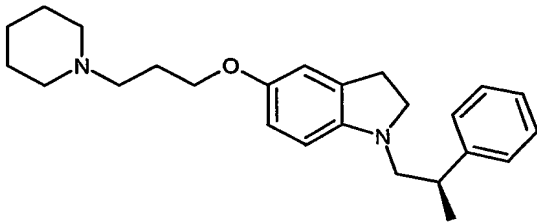
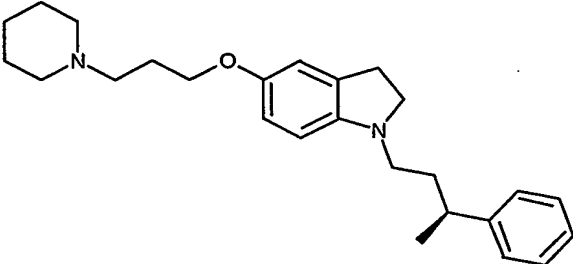
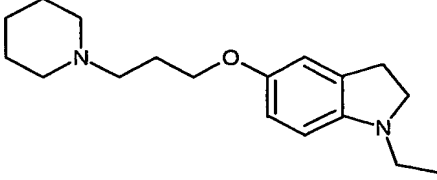
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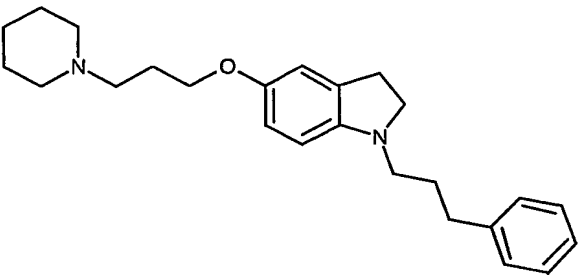
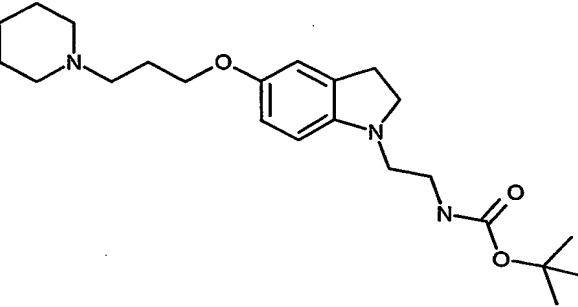
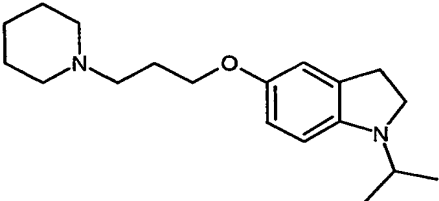
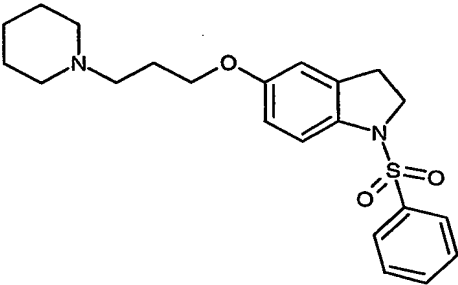
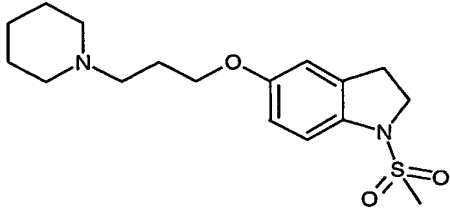
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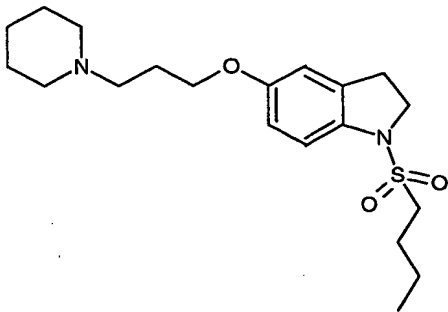
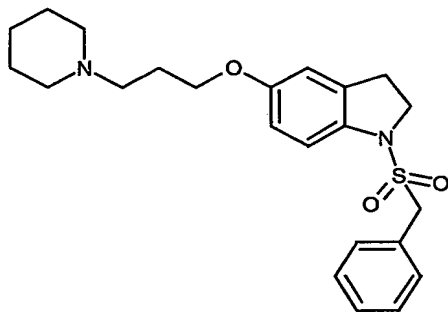
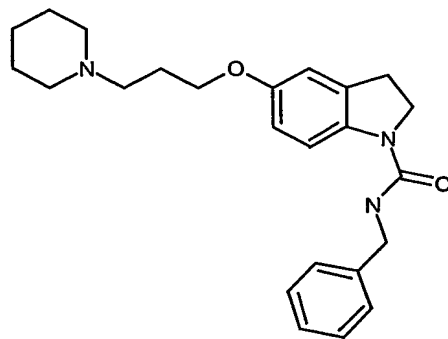
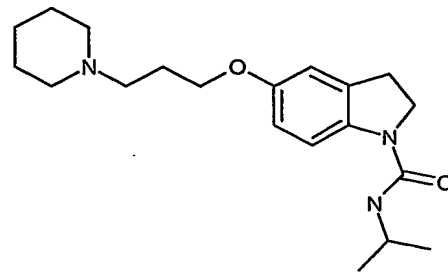
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	22	 <chem>CNCC1CCN(C1)CCOC2=CC=C3C(=C2)N(CCC4=CC=CC=C4N5C=CC=CC=C5)C3</chem>	
	23	 <chem>c1ccccc1CN2C=CC=CC=C2C3=CC=C4C(=C3)N(CCC5CCN(CC5)CC6=CC=CC=C6O4)C</chem>	
	24	 <chem>ClC1=CC(=C(C(=C1)N2C=CC=CC=C2C3=CC=C4C(=C3)N(CCC5CCN(CC5)CC6=CC=CC=C6O4)C)C)C</chem>	
	25	 <chem>c1cc(s1)CN2C=CC=CC=C2C3=CC=C4C(=C3)N(CCC5CCN(CC5)CC6=CC=CC=C6O4)C</chem>	
	26	 <chem>c1cc(s1)CN2C=CC=CC=C2C3=CC=C4C(=C3)N(CCC5CCN(CC5)CC6=CC=CC=C6O4)C</chem>	

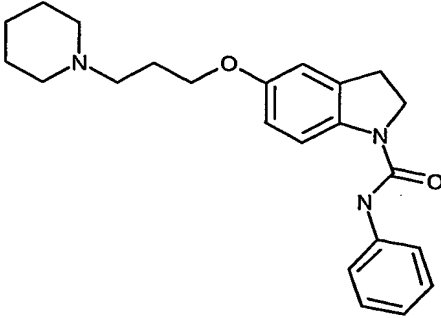
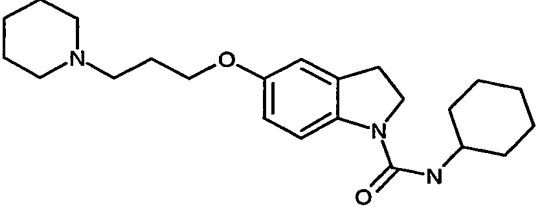
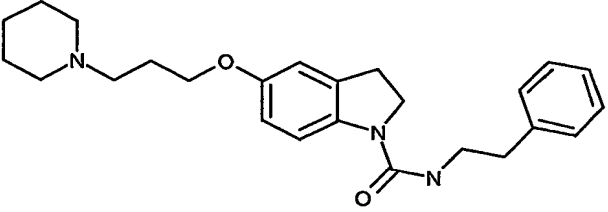
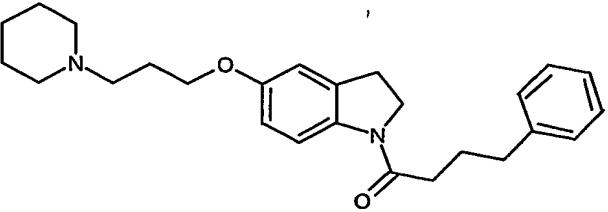
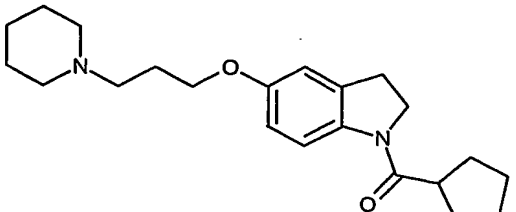
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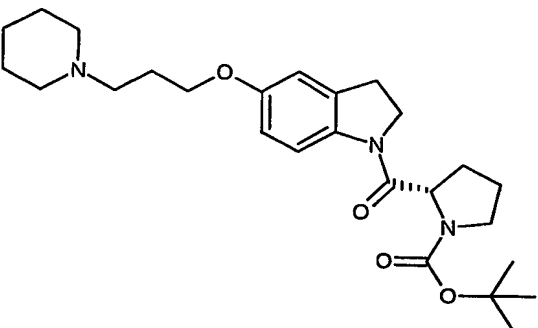
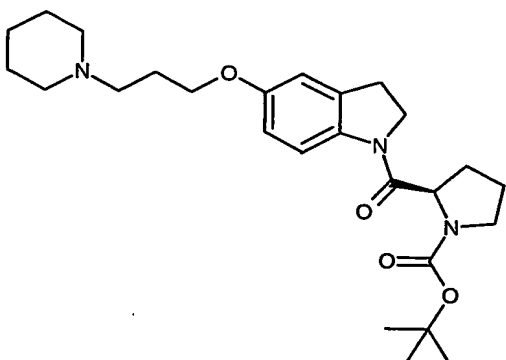
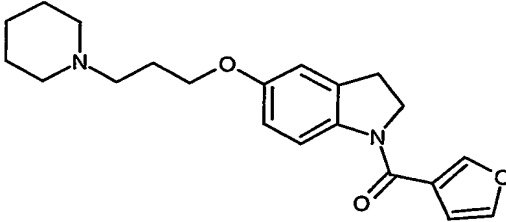
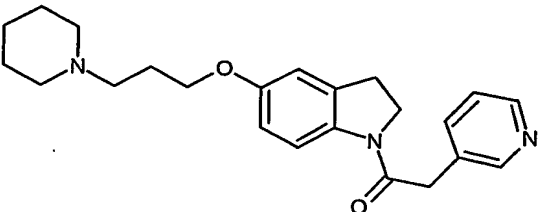
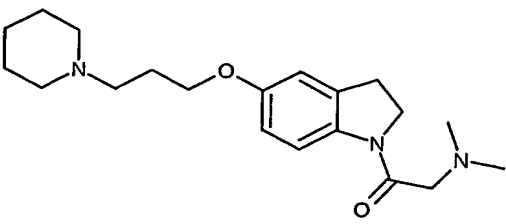


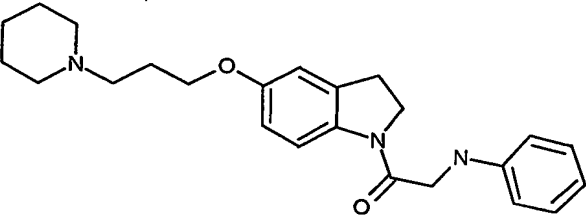
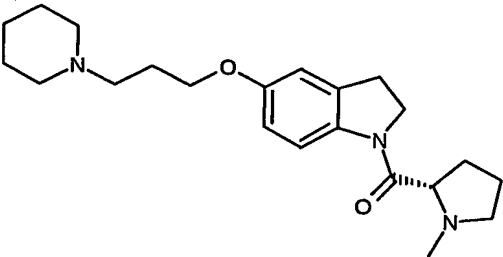
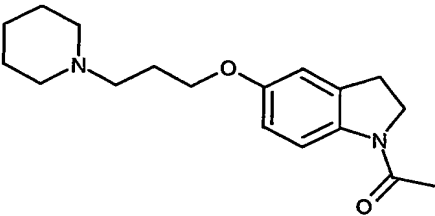
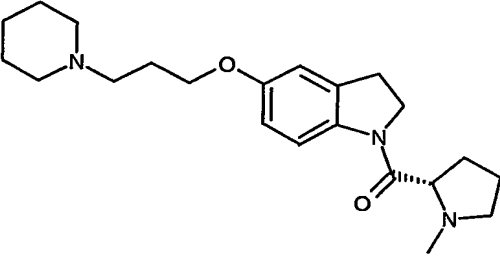
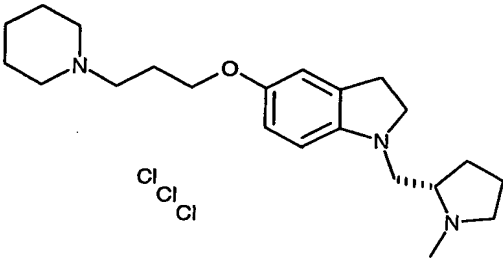
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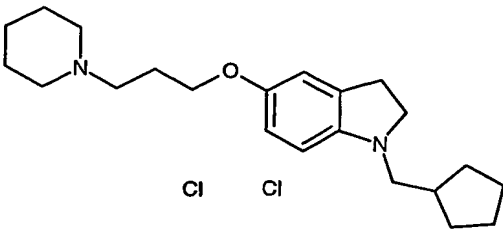
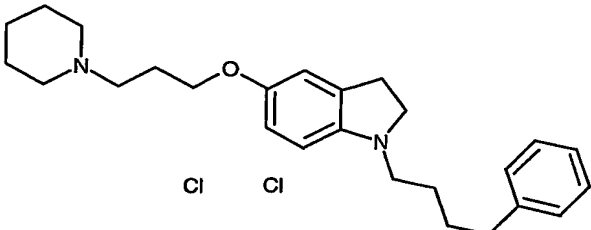
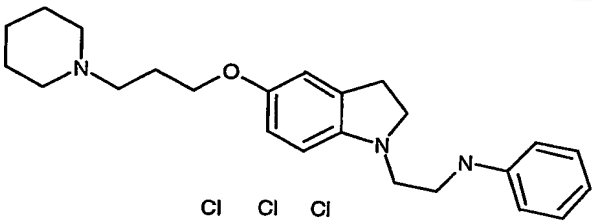
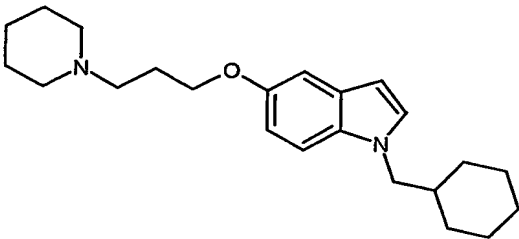
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	51	 <p>Chiral</p>	
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	57	 Chiral	
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	59	 Chiral	
	60	 Chiral	

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or a pharmaceutically acceptable salt or solvate thereof.

8. A pharmaceutical composition which comprises a compound of any of claims 1-7 and a pharmaceutically acceptable carrier.
- 5 9. A method of selectively increasing histamine levels in cells by contacting the cells with an antagonist of the histamine H<sub>3</sub> receptor, said antagonists comprising a compound of any of claims 1-7.
10. The method of Claim 9 wherein the antagonist is a pharmaceutical composition of claim 8.
- 10 11. The method of Claim 9 wherein the antagonist is characterized by selectively binding the histamine H<sub>3</sub>R receptor, relative to the histamine H<sub>4</sub>R receptor.

12. A method for treatment or prevention of obesity which comprises administering to a subject in need of such treatment or prevention an effective amount of a compound of any of Claims 1-7.
13. The method of Claim 12 wherein the antagonist is a pharmaceutical composition  
5 of claim 8.
14. A method for treatment or prevention of a disorder or disease in which inhibition of the histamine H3 receptor has a beneficial effect which comprises administering to a subject in need of such treatment or prevention an effective amount of a compound of any of claims 1-7.